

Connecting via Winsock to STN

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LOGINID:SSPTAEAL1624

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

NEWS	1		Web Page for STN Seminar Schedule - N. America
NEWS	2	JUL 02	LMEDLINE coverage updated
NEWS	3	JUL 02	SCISEARCH enhanced with complete author names
NEWS	4	JUL 02	CHEMCATS accession numbers revised
NEWS	5	JUL 02	CA/CAPplus enhanced with utility model patents from China
NEWS	6	JUL 16	CAPplus enhanced with French and German abstracts
NEWS	7	JUL 18	CA/CAPplus patent coverage enhanced
NEWS	8	JUL 26	USPATFULL/USPAT2 enhanced with IPC reclassification
NEWS	9	JUL 30	USGENE now available on STN
NEWS	10	AUG 06	CAS REGISTRY enhanced with new experimental property tags
NEWS	11	AUG 06	BEILSTEIN updated with new compounds
NEWS	12	AUG 06	FSTA enhanced with new thesaurus edition
NEWS	13	AUG 13	CA/CAPplus enhanced with additional kind codes for granted patents
NEWS	14	AUG 20	CA/CAPplus enhanced with CAS indexing in pre-1907 records
NEWS	15	AUG 27	Full-text patent databases enhanced with predefined patent family display formats from INPADOCDB
NEWS	16	AUG 27	USPATOLD now available on STN
NEWS	17	AUG 28	CAS REGISTRY enhanced with additional experimental spectral property data
NEWS	18	SEP 07	STN AnaVist, Version 2.0, now available with Derwent World Patents Index
NEWS	19	SEP 13	FORIS renamed to SOFIS
NEWS	20	SEP 13	INPADOCDB enhanced with monthly SDI frequency
NEWS	21	SEP 17	CA/CAPplus enhanced with printed CA page images from 1967-1998
NEWS	22	SEP 17	CAPplus coverage extended to include traditional medicine patents
NEWS	23	SEP 24	EMBASE, EMBAL, and LEMBASE reloaded with enhancements
NEWS EXPRESS	19	SEPTEMBER 2007:	CURRENT WINDOWS VERSION IS V8.2, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 19 SEPTEMBER 2007.
NEWS HOURS			STN Operating Hours Plus Help Desk Availability
NEWS LOGIN			Welcome Banner and News Items
NEWS IPC8			For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that specific topic.

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* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 16:41:51 ON 24 SEP 2007

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'REGISTRY' ENTERED AT 16:41:58 ON 24 SEP 2007

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STRUCTURE FILE UPDATES: 23 SEP 2007 HIGHEST RN 947726-74-1

DICTIONARY FILE UPDATES: 23 SEP 2007 HIGHEST RN 947726-74-1

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 29, 2007

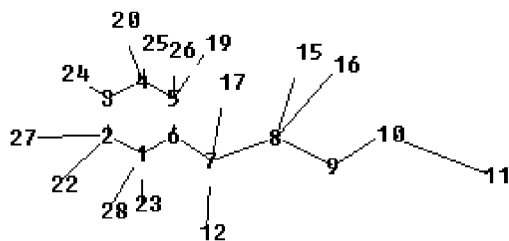
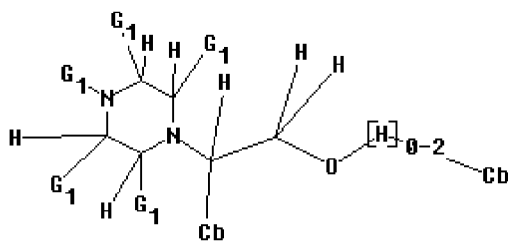
Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stdoc/properties.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10599824.str



```

chain nodes :
7 8 9 10 11 12 15 16 17 19 20 22 23 24 25 26 27 28
ring nodes :
1 2 3 4 5 6
chain bonds :
1-23 1-28 2-22 2-27 3-24 4-20 4-25 5-19 5-26 6-7 7-8 7-12 7-17 8-9
8-15 8-16 9-10 10-11
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6
exact/norm bonds :
1-2 1-6 1-23 2-3 2-22 3-4 3-24 4-5 4-20 5-6 5-19 6-7 8-9
exact bonds :
1-28 2-27 4-25 5-26 7-8 7-12 7-17 8-15 8-16 9-10 10-11
isolated ring systems :
containing 1 :

```

G1:H,X,Ak,O

```

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS
11:Atom 12:Atom 15:CLASS 16:CLASS 17:CLASS 19:CLASS 20:CLASS 22:CLASS
23:CLASS 24:CLASS
25:CLASS 26:CLASS 27:CLASS 28:CLASS

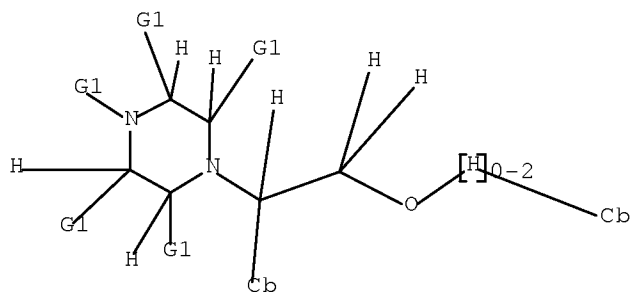
```

L1 STRUCTURE UPLOADED

=> d 11

L1 HAS NO ANSWERS

L1 STR



G1 H, X, Ak, O

Structure attributes must be viewed using STN Express query preparation.

=> s l1 full

FULL SEARCH INITIATED 16:42:20 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 258294 TO ITERATE

100.0% PROCESSED 258294 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.04

L2 1 SEA SSS FUL L1

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

172.10

172.31

FILE 'CAPLUS' ENTERED AT 16:42:31 ON 24 SEP 2007

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FILE COVERS 1907 - 24 Sep 2007 VOL 147 ISS 14

FILE LAST UPDATED: 23 Sep 2007 (20070923/ED)

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<http://www.cas.org/infopolicy.html>

=> s l1 full

REGISTRY INITIATED

Substance data SEARCH and crossover from CAS REGISTRY in progress...
Use DISPLAY HITSTR (or FHITSTR) to directly view retrieved structures.

FULL SEARCH INITIATED 16:42:39 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 258294 TO ITERATE

100.0% PROCESSED 258294 ITERATIONS 1 ANSWERS
SEARCH TIME: 00.00.04

L3 1 SEA SSS FUL L1

L4 1 L3

=> file caplus
COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
0.47	345.35

FULL ESTIMATED COST

FILE 'CAPLUS' ENTERED AT 16:42:49 ON 24 SEP 2007
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FILE LAST UPDATED: 23 Sep 2007 (20070923/ED)

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=> s l2 full

L5 1 L2

=> d ibib abs

L5 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2007 ACS on STN
ACCESSION NUMBER: 2004:589533 CAPLUS Full-text
DOCUMENT NUMBER: 141:140464
TITLE: N-(substituted arylmethyl)-4-(disubstituted methyl)piperidines and piperazines
INVENTOR(S): Ding, Ping; Henrie, Robert N., II; Cohen, Daniel H.; Lyga, John W.; Rosen, David S.; Theodoridis, George;

Zhang, Qun; Yeager, Walter H.; Donovan, Stephen F.;
 Zhang, Steven Shunxiang; Shulman, Inna; Yu, Seong Jae;
 Wnag, Gouzhi; Zhang, Y. Larry; Gopalsamy, Ariamala;
 Warkentin, Dennis L.; Rensner, Paul E.; Silverman, Ian
 R.; Cullen, Thomas G.

PATENT ASSIGNEE(S): FMC Corporation, USA
 SOURCE: PCT Int. Appl., 105 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

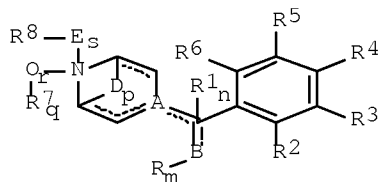
LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

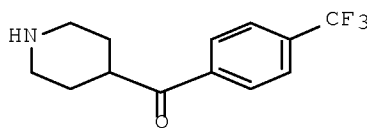
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004060865	A2	20040722	WO 2003-US39046	20031208
WO 2004060865	A3	20041104		
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
RW:	BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
AU 2003296373	A1	20040729	AU 2003-296373	20031208
EP 1572668	A2	20050914	EP 2003-814673	20031208
R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK			
BR 2003016747	A	20051018	BR 2003-16747	20031208
CN 1729178	A	20060201	CN 2003-80106750	20031208
CN 1744895	A	20060308	CN 2003-80109445	20031208
JP 2006511621	T	20060406	JP 2005-508564	20031208
IN 2005DN02489	A	20061229	IN 2005-DN2489	20050609
IN 2005DN02485	A	20070427	IN 2005-DN2485	20050609
ZA 2005004870	A	20060426	ZA 2005-4870	20050614
ZA 2005004871	A	20060426	ZA 2005-4871	20050614
MX 2005PA06426	A	20050908	MX 2005-PA6426	20050615
US 2006166962	A1	20060727	US 2006-538997	20060208
PRIORITY APPLN. INFO.:			US 2002-434718P	P 20021218
			US 2003-495059P	P 20030814
			WO 2003-US39046	W 20031208

OTHER SOURCE(S): MARPAT 141:140464
 GI



I



II

AB Title compds. I [m, n, q, r, s = 0-1; p = 0-3; A = CH, N forming a 6-membered azine ring selected from piperidine or piperazine; R2-6 = H, halo, alkyl, etc.; B = O; with provisions] are prepared For instance, 4-bromobenzotrifluoride is transmetalated (THF, n-BuLi, -75°) and treated with tert-Bu 4-[N-methoxy-N-methylcarbamoyl]piperidine-1- carboxylate to give tert-Bu 4-[(4-(trifluoromethyl)phenyl)carbonyl]piperid ine-1-carboxylate. This intermediate is deprotected to give II. II gave 100% mortality and 100% growth inhibition of tobacco budworms.

=> d hitstr

L5 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2007 ACS on STN

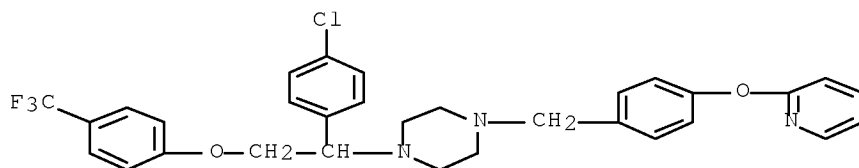
IT 725231-94-7P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(N-(substituted arylmethyl)-4-(disubstituted methyl)piperidines and piperazines)

RN 725231-94-7 CAPLUS

CN Piperazine, 1-[1-(4-chlorophenyl)-2-[4-(trifluoromethyl)phenoxy]ethyl]-4-[[4-(2-pyridinyloxy)phenyl]methyl]- (9CI) (CA INDEX NAME)



=> file reg

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION

FULL ESTIMATED COST

8.56	353.91
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DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION

CA SUBSCRIBER PRICE

-0.78	-0.78
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FILE 'REGISTRY' ENTERED AT 16:47:02 ON 24 SEP 2007

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STRUCTURE FILE UPDATES: 23 SEP 2007 HIGHEST RN 947726-74-1

DICTIONARY FILE UPDATES: 23 SEP 2007 HIGHEST RN 947726-74-1

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TSCA INFORMATION NOW CURRENT THROUGH June 29, 2007

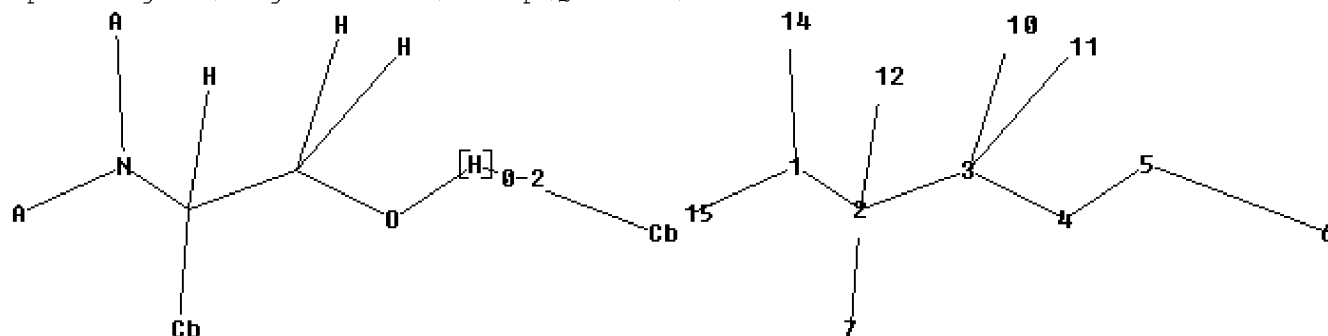
Please note that search-term pricing does apply when
conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10599824broad.str



chain nodes :

2 3 4 5 6 7 10 11 12

ring/chain nodes :

1 14 15

chain bonds :

1-2 1-14 1-15 2-3 2-7 2-12 3-4 3-10 3-11 4-5 5-6

exact/norm bonds :

1-2 1-14 1-15 3-4

exact bonds :

2-3 2-7 2-12 3-10 3-11 4-5 5-6

G1:H,X,Ak,O

Match level :

1:Atom 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:Atom 7:Atom 10:CLASS 11:CLASS

12:CLASS

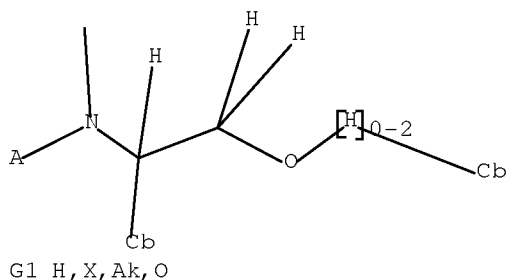
14:CLASS 15:CLASS

L6 STRUCTURE UPLOADED

=> d 16

L6 HAS NO ANSWERS

L6 STR



Structure attributes must be viewed using STN Express query preparation.

=> s 16 full

FULL SEARCH INITIATED 16:50:46 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 6037621 TO ITERATE

8.6% PROCESSED 516823 ITERATIONS 7 ANSWERS

15.8% PROCESSED 955056 ITERATIONS 7 ANSWERS

16.6% PROCESSED 1000000 ITERATIONS 7 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.39

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
BATCH **INCOMPLETE**
PROJECTED ITERATIONS: 6037621 TO 6037621
PROJECTED ANSWERS: 23 TO 61

L7 7 SEA SSS FUL L6

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	175.25	529.16
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-0.78

FILE 'CAPLUS' ENTERED AT 16:51:32 ON 24 SEP 2007

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FILE COVERS 1907 - 24 Sep 2007 VOL 147 ISS 14
FILE LAST UPDATED: 23 Sep 2007 (20070923/ED)

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<http://www.cas.org/infopolicy.html>

=> s 17 full

L8 2 L7

=> d ibib abs hitstr tot

L8 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 2006:908614 CAPLUS Full-text

DOCUMENT NUMBER: 145:454704

TITLE: Effect of the Phosphoryl Substituent in the Linear
Nitron on the Spin Trapping of Superoxide Radical and
the Stability of the Superoxide Adduct: Combined
Experimental and Theoretical Studies

AUTHOR(S): Liu, Yang-Ping; Wang, Lan-Fen; Nie, Zhou; Ji,
Yi-Qiong; Liu, Yang; Liu, Ke-Jian; Tian, Qiu

CORPORATE SOURCE: State Key Laboratory for Structural Chemistry of
Unstable and Stable Species, Center for Molecular
Science, Institute of Chemistry, Chinese Academy of
Sciences, Beijing, 100080, Peop. Rep. China

SOURCE: Journal of Organic Chemistry (2006), 71(20), 7753-7762
CODEN: JOCEAH; ISSN: 0022-3263

PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 145:454704

AB A new phosphorylated linear nitron N-(4-hydroxybenzylidene)-1-
diethoxyphosphoryl-1-methylethylamine N-oxide (4-HOPPN) was synthesized, and
its X-ray structure was determined. The spin trapping ability of various kinds
of free radicals by 4-HOPPN was evaluated. Kinetic study of decay of the
superoxide spin adduct (4-HOPPN-OOH) shows the half-life time of 8.8 min. On
the basis of the X-ray structural coordinates, theor. analyses using d.
functional theory (DFT) calcns. at the B3LYP/6-31+G(d,p)//B3LYP/6-31G(d) level
were performed on spin-trapping reactions of superoxide radical with 4-HOPPN
and PBN and three possible decay routes for their corresponding superoxide
adducts. The comparative calcns. on the spin-trapping reactions with
superoxide radical predicted that both spin traps share an identical reaction
type and have comparable potency when spin trapping superoxide radical. Anal.
of the optimized geometries of 4-HOPPN-OOH and PBN-OOH reveals that an
introduction of the phosphoryl group can efficiently stabilize the spin adduct
through the intramol. H-bonds, the intramol. nonbonding attractive
interactions, as well as the bulky steric protection. Examination of the
decomposition thermodyn. of 4-HOPPN-OOH and PBN-OOH further supports the
stabilizing role of the phosphoryl group to a linear phosphorylated spin
adduct.

IT 913260-59-0

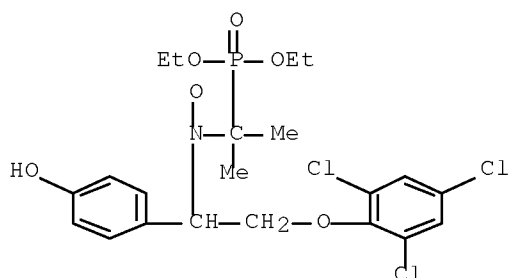
RL: FMU (Formation, unclassified); PRP (Properties); FORM (Formation,
nonpreparative)

(effect of phosphoryl substituent in linear nitron on spin trapping of
superoxide radical and stability of superoxide)

RN 913260-59-0 CAPLUS

CN Nitroxide, 1-(diethoxyphosphinyl)-1-methylethyl 1-(4-hydroxyphenyl)-2-

(2,4,6-trichlorophenoxy)ethyl (9CI) (CA INDEX NAME)



REFERENCE COUNT: 73 THERE ARE 73 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L8 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 2006:874491 CAPLUS Full-text

DOCUMENT NUMBER: 145:471200

TITLE: Synthesis and evaluation of anilinohexafluoroisopropanols as activators/modulators of LXR α and β

AUTHOR(S): Panday, Narendra; Benz, Jorg; Blum-Kaelin, Denise; Bourgeaux, Vanessa; Dehmlow, Henrietta; Hartman, Peter; Kuhn, Bernd; Ratni, Hassen; Warot, Xavier; Wright, Matthew B.

CORPORATE SOURCE: Pharmaceuticals Division, Preclinical Research, F. Hoffmann-La Roche Ltd., Basel, CH-4070, Switz.

SOURCE: Bioorganic & Medicinal Chemistry Letters (2006), 16(19), 5231-5237
CODEN: BMCLE8; ISSN: 0960-894X

PUBLISHER: Elsevier Ltd.

DOCUMENT TYPE: Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 145:471200

AB A series of branched and unbranched anilinohexafluoroisopropanols related to the known sulfonamide T0901317 were prepared and evaluated as activators/modulators of both LXR α and LXR β . A structure-activity relationship was established and compds. with high potency on both the receptors were identified. Many compds. showed a tendency toward selectivity for LXR β vs. LXR α . Several analogs were evaluated for effects on plasma lipoprotein levels in mice. A few of these significantly raised HDL-cholesterol levels in plasma but showed markedly different effects on liver triglyceride content, suggesting that this series may yield candidates with improved efficacy/safety profiles compared to existing mols.

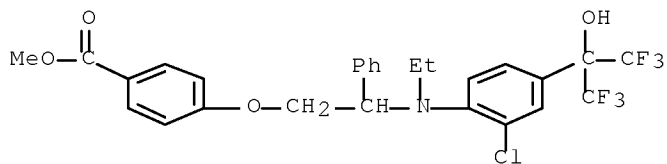
IT 913619-59-7F 913619-60-0F 913619-61-1F

RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent)

(synthesis and evaluation of anilinohexafluoroisopropanols as activators/modulators of LXR α and β)

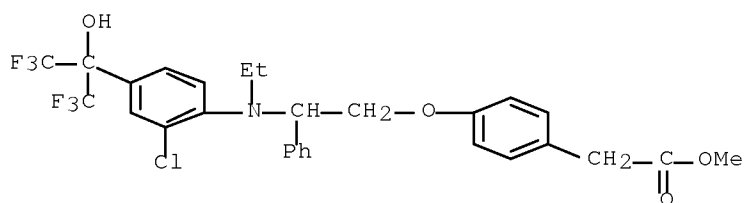
RN 913619-59-7 CAPLUS

CN Benzoic acid, 4-[2-[[[2-chloro-4-[2,2,2-trifluoro-1-hydroxy-1-(trifluoromethyl)ethyl]phenyl]ethylamino]-2-phenylethoxy]-, methyl ester (CA INDEX NAME)



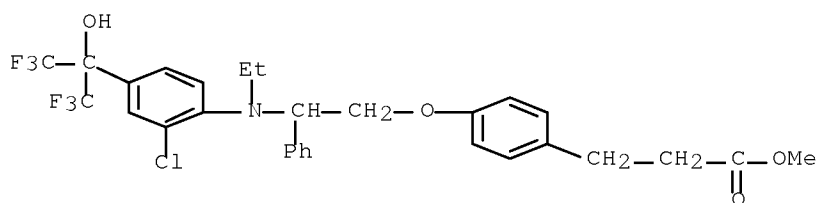
RN 913619-60-0 CAPLUS

CN Benzenepropanoic acid, 4-[2-[[2-chloro-4-[2,2,2-trifluoro-1-hydroxy-1-(trifluoromethyl)ethyl]phenyl]ethylamino]-2-phenylethoxy]-, methyl ester (CA INDEX NAME)



RN 913619-61-1 CAPLUS

CN Benzenepropanoic acid, 4-[2-[[2-chloro-4-[2,2,2-trifluoro-1-hydroxy-1-(trifluoromethyl)ethyl]phenyl]ethylamino]-2-phenylethoxy]-, methyl ester (CA INDEX NAME)

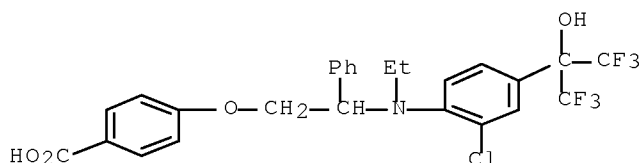


IT 913619-62-2P 913619-63-3P 913619-64-4P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)
(synthesis and evaluation of anilinohexafluoroisopropanols as activators/modulators of LXR α and β)

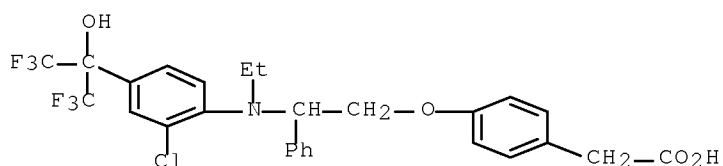
RN 913619-62-2 CAPLUS

CN Benzoic acid, 4-[2-[[2-chloro-4-[2,2,2-trifluoro-1-hydroxy-1-(trifluoromethyl)ethyl]phenyl]ethylamino]-2-phenylethoxy]- (CA INDEX NAME)



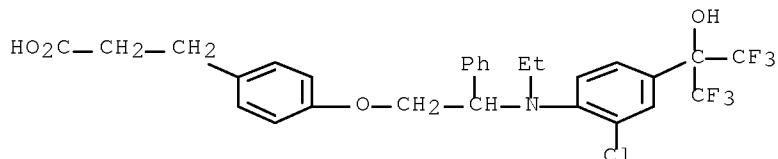
RN 913619-63-3 CAPLUS

CN Benzenepropanoic acid, 4-[2-[[2-chloro-4-[2,2,2-trifluoro-1-hydroxy-1-(trifluoromethyl)ethyl]phenyl]ethylamino]-2-phenylethoxy]- (CA INDEX NAME)



RN 913619-64-4 CAPLUS

CN Benzenepropanoic acid, 4-[2-[[2-chloro-4-[2,2,2-trifluoro-1-hydroxy-1-(trifluoromethyl)ethyl]phenyl]ethylamino]-2-phenylethoxy]- (CA INDEX NAME)



REFERENCE COUNT: 24 THERE ARE 24 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> file reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

12.89

542.05

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

CA SUBSCRIBER PRICE

-1.56

-2.34

FILE 'REGISTRY' ENTERED AT 16:54:45 ON 24 SEP 2007

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 23 SEP 2007 HIGHEST RN 947726-74-1
DICTIONARY FILE UPDATES: 23 SEP 2007 HIGHEST RN 947726-74-1

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 29, 2007

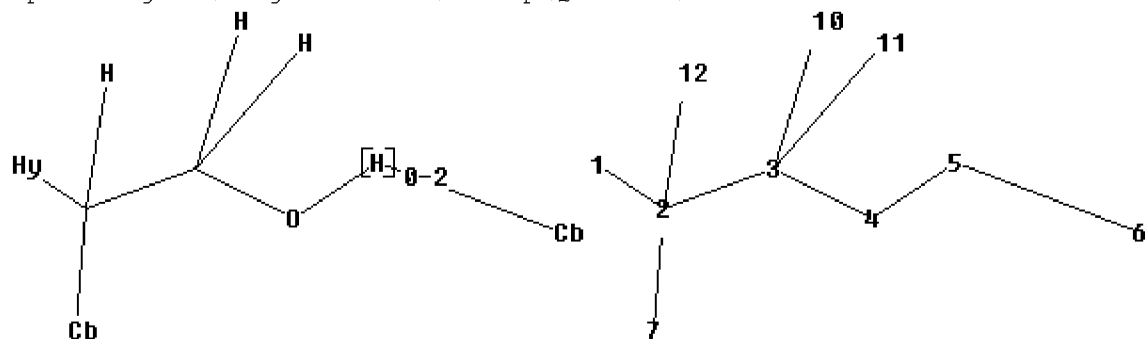
Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stdoc/properties.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10599824last.str



chain nodes :

1 2 3 4 5 6 7 10 11 12

chain bonds :

1-2 2-3 2-7 2-12 3-4 3-10 3-11 4-5 5-6

exact/norm bonds :

1-2 3-4

exact bonds :

2-3 2-7 2-12 3-10 3-11 4-5 5-6

G1:H,X,Ak,O

Match level :

1:Atom 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:Atom 7:Atom 10:CLASS 11:CLASS

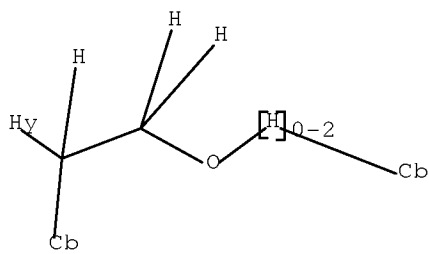
12:CLASS

L9 STRUCTURE UPLOADED

=> d 19

L9 HAS NO ANSWERS

STR



Structure attributes must be viewed using STN Express query preparation.

```
=> s 19 full
FULL SEARCH INITIATED 16:55:06 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 25530888 TO ITERATE
```

```

0.8% PROCESSED      200310 ITERATIONS                      0 ANSWERS
1.6% PROCESSED      398021 ITERATIONS                      0 ANSWERS
2.4% PROCESSED      605940 ITERATIONS                      0 ANSWERS
3.5% PROCESSED      890366 ITERATIONS                      0 ANSWERS
3.7% PROCESSED      932709 ITERATIONS                      0 ANSWERS
3.9% PROCESSED     1000000 ITERATIONS                      0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.01.38

```

```

FULL FILE PROJECTIONS:  ONLINE  **INCOMPLETE**
                        BATCH    **INCOMPLETE**
PROJECTED ITERATIONS:   25530888 TO 25530888
PROJECTED ANSWERS:      0 TO      0

```

L10 0 SEA SSS FUL L9

=> d his

(FILE 'HOME' ENTERED AT 16:41:51 ON 24 SEP 2007)

FILE 'REGISTRY' ENTERED AT 16:41:58 ON 24 SEP 2007

L1 STRUCTURE UPLOADED

L2 1 S L1 FULL

FILE 'CAPLUS' ENTERED AT 16:42:31 ON 24 SEP 2007

S L1

FILE 'REGISTRY' ENTERED AT 16:42:39 ON 24 SEP 2007

L3 1 S L1 FULL

FILE 'CAPLUS' ENTERED AT 16:42:43 ON 24 SEP 2007

L4 1 S L3 FULL

FILE 'CAPLUS' ENTERED AT 16:42:49 ON 24 SEP 2007

L5 1 S L2 FULL

FILE 'REGISTRY' ENTERED AT 16:47:02 ON 24 SEP 2007

L6 STRUCTURE UPLOADED

L7 7 S L6 FULL

FILE 'CAPLUS' ENTERED AT 16:51:32 ON 24 SEP 2007

L8 2 S L7 FULL

FILE 'REGISTRY' ENTERED AT 16:54:45 ON 24 SEP 2007

L9 STRUCTURE UPLOADED

L10 0 S L9 FULL

=> log yu

'YU' IS NOT VALID HERE

For an explanation, enter "HELP LOGOFF".

=> log y

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

173.45

715.50

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

CA SUBSCRIBER PRICE

0.00

-2.34

STN INTERNATIONAL LOGOFF AT 16:57:12 ON 24 SEP 2007